

# An efficient and spectrally accurate numerical method for computing dynamics of rotating Bose-Einstein condensates

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## Abstract

In this paper, we propose an efficient and spectrally accurate numerical method for computing the dynamics of rotating Bose-Einstein condensates (BEC) in two dimensions (2D) and 3D based on the Gross-Pitaevskii equation (GPE) with an angular momentum rotation term. By applying a time-splitting technique for decoupling the nonlinearity and properly using the alternating direction implicit (ADI) technique for the coupling in the angular momentum rotation term in the GPE, at every time step, the GPE in rotational frame is decoupled into a nonlinear ordinary differential equation (ODE) and two partial differential equations with constant coefficients. This allows us to develop new time-splitting spectral (TSSP) methods for computing the dynamics of BEC in a rotational frame. The new numerical method is explicit, unconditionally stable, and of spectral accuracy in space and second order accuracy in time. Moreover, it is time reversible and time transverse invariant, and conserves the position density in the discretized level if the GPE does. Extensive numerical results are presented to confirm the above properties of the new numerical method for rotating BEC in 2D & 3D.

*Key Words:* Rotating Bose-Einstein condensates, Gross-Pitaevskii equation, angular momentum rotation, time spitting.

## 1 Introduction

Since the first experimental creation of a quantized vortex in a gaseous Bose-Einstein condensate (BEC) [29, 1, 30, 35], there has been significantly experimental and theoretical advances in the field of research [3, 34, 2, 11, 5, 13, 23, 14, 17, 19, 22].

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Several experimental methods of vortex creation are currently in use, including phase imprinting [30, 40], cooling of a rotating normal gas [24], and conversion of spin angular momentum into orbital angular momentum by reversal of the magnetic bias field in an Ioffe-Pritchard trap [26, 27, 33]. The topic of this paper is to propose an efficient and spectrally accurate numerical method for studying quantized vortex dynamics in a BEC by imposing a laser beam rotating with an angular velocity on the magnetic trap holding the atoms to create a harmonic anisotropic potential.

At temperatures  $T$  much smaller than the critical condensation temperature  $T_c$ , under mean field theory, the properties of a BEC in a rotational frame are modelled by the well-known time-dependent Gross-Pitaevskii equation (GPE) with an angular momentum rotation term [3, 14, 22]:

$$i\hbar \frac{\partial \psi(\mathbf{x}, t)}{\partial t} = \left( -\frac{\hbar^2}{2m} \nabla^2 + V(\mathbf{x}) + NU_0 |\psi|^2 - \Omega L_z \right) \psi(\mathbf{x}, t), \quad \mathbf{x} \in \mathbb{R}^3, \quad t \geq 0, \quad (1.1)$$

where  $\mathbf{x} = (x, y, z)^T$  is the Cartesian coordinate vector,  $\psi(\mathbf{x}, t)$  is the complex-valued macroscopic wave function,  $m$  is the atomic mass,  $\hbar$  is the Planck constant,  $N$  is the number of atoms in the condensate,  $\Omega$  is the angular velocity of rotating laser beam,  $V(\mathbf{x}) = \frac{m}{2} (\omega_x^2 x^2 + \omega_y^2 y^2 + \omega_z^2 z^2)$  with  $\omega_x, \omega_y$  and  $\omega_z$  being the trap frequencies in  $x$ -,  $y$ - and  $z$ -direction respectively.  $U_0 = \frac{4\pi\hbar^2 a_s}{m}$  describes the interaction between atoms in the condensate with  $a_s$  the  $s$ -wave scattering length, and  $L_z$  is the  $z$ -component of the angular momentum. It is convenient to normalize the wave function by requiring

$$\|\psi(\cdot, t)\|^2 := \int_{\mathbb{R}^3} |\psi(\mathbf{x}, t)|^2 d\mathbf{x} = 1. \quad (1.2)$$

After proper nondimensionalization and dimension reduction, we can obtain the following dimensionless GPE with an angular momentum rotation term in the  $d$ -dimensions ( $d = 2, 3$ ) [13, 5, 41]:

$$i \frac{\partial \psi(\mathbf{x}, t)}{\partial t} = -\frac{1}{2} \nabla^2 \psi + V_d(\mathbf{x}) \psi + \beta_d |\psi|^2 \psi - \Omega L_z \psi, \quad \mathbf{x} \in \mathbb{R}^d, \quad t \geq 0, \quad (1.3)$$

$$\psi(\mathbf{x}, 0) = \psi_0(\mathbf{x}), \quad \mathbf{x} \in \mathbb{R}^d, \quad \text{with} \quad \|\psi_0\|^2 := \int_{\mathbb{R}^d} |\psi_0(\mathbf{x})|^2 d\mathbf{x} = 1, \quad (1.4)$$

where  $L_z = -i(x\partial_y - y\partial_x)$  and

$$V_d(\mathbf{x}) = \begin{cases} (\gamma_x^2 x^2 + \gamma_y^2 y^2) / 2, & d = 2, \\ (\gamma_x^2 x^2 + \gamma_y^2 y^2 + \gamma_z^2 z^2) / 2, & d = 3; \end{cases} \quad (1.5)$$

with  $\gamma_x, \gamma_y$  and  $\gamma_z$  being constants.

In order to study effectively the dynamics of BEC, especially in the strong repulsive interaction regime, i.e.  $\beta_d \gg 1$  in (1.3), an efficient and accurate numerical method is one of the key issues. For non-rotating BEC, i.e.  $\Omega = 0$  in (1.3), many efficient and spectrally accurate numerical methods were proposed in the literatures [7, 6, 11, 4, 12], and they were demonstrated that they are much better than the

low-order finite difference methods [15, 32, 31, 16]. Thus they were applied to study collapse and explosion of BEC in 3D [8] and multi-component BEC [4] which are the very challenging problems in numerical simulation of BEC. Due to the appearance of the angular momentum rotation term in the GPE (1.3), new numerical difficulties are introduced. Currently, the numerical methods used in the physics literature for studying dynamics of rotating BEC remain limited [25, 39], and they usually are low-order finite difference methods. Recently, some efficient and accurate numerical methods were designed for computing dynamics of rotating BEC. For example, Bao, Du and Zhang [5] proposed a numerical method for computing dynamics of rotating BEC by applying a time-splitting technique for decoupling the nonlinearity in the GPE and adopting the polar coordinates or cylindrical coordinates so as to make the coefficient of the angular momentum rotation term constant. The method is time reversible, unconditionally stable, implicit but can be solved very efficiently, and conserves the total density. It is of spectral accuracy in transverse direction, but usually of second or fourth-order accuracy in radial direction. Another numerical method is the leap-frog spectral method used for studying vortex lattice dynamics in rotating BEC [41]. This method is explicit, time reversible, of spectral accuracy in space and second order accuracy in time. But it has a stability constraint for time step [41]. The aim of this paper is to develop a numerical method which enjoys advantages of both the above two numerical methods. That is to say, the method is explicit, unconditionally stable, time reversible, time transverse invariant, and of spectral accuracy in space. We shall present such an efficient, unconditionally stable and accurate numerical method for discretizing the GPE in a rotational frame by applying a time-splitting technique and an ADI technique, and constructing appropriately spectral basis functions.

The paper is organized as follows. In section 2, we review some properties of GPE in a rotational frame (1.3) including conservation laws and analytical solutions of condensate widths. In section 3, we present a new time-splitting Fourier pseudospectral method for efficient and accurate simulation of GPE (1.3) in 2D & 3D. In section 4, extensive numerical results are reported to demonstrate the efficiency and spectral resolution in space of our new numerical method. Finally some conclusions are drawn in section 5.

## 2 Some properties of the GPE

For the convenience of the reader, in this section, we will review some properties of the GPE with an angular momentum rotation term (1.3), which will be used to test our new numerical method proposed in the next section. First of all, the GPE (1.3) is time reversible and time transverse invariant. Second, it has at least two important invariants which are the *normalization of the wave function*

$$N(\psi) = \int_{\mathbb{R}^d} |\psi(\mathbf{x}, t)|^2 d\mathbf{x} \equiv \int_{\mathbb{R}^d} |\psi(\mathbf{x}, 0)|^2 d\mathbf{x} = N(\psi_0) = 1, \quad t \geq 0, \quad (2.1)$$

and the energy

$$\begin{aligned} E_{\beta,\Omega}(\psi) &= \int_{\mathbb{R}^d} \left[ \frac{1}{2} |\nabla \psi|^2 + V_d(\mathbf{x}) |\psi|^2 + \frac{\beta_d}{2} |\psi|^4 - \Omega \psi^* L_z \psi \right] d\mathbf{x} \\ &= E_{\beta_d,\Omega}(\psi_0), \quad t \geq 0, \end{aligned} \quad (2.2)$$

where  $f^*$  and  $\text{Im}(f)$  denote the conjugate and the imaginary part of the function  $f$  respectively. Third, it was proven that at least for radial symmetric trap in 2D or cylindrical symmetric trap in 3D, i.e.,  $\gamma_x = \gamma_y$  in (1.3), the angular momentum expectation and energy for non-rotating part are conserved [5], that is, for any given initial data  $\psi_0(\mathbf{x})$  in (1.4),

$$\langle L_z \rangle(t) \equiv \langle L_z \rangle(0), \quad E_{\beta,0}(\psi) \equiv E_{\beta,0}(\psi_0), \quad t \geq 0, \quad (2.3)$$

where the angular momentum expectation which is a measure of the vortex flux is defined as

$$\langle L_z \rangle(t) := \int_{\mathbb{R}^d} \psi^*(\mathbf{x}, t) L_z \psi(\mathbf{x}, t) d\mathbf{x} = i \int_{\mathbb{R}^d} \psi^*(\mathbf{x}, t) (y \partial_x - x \partial_y) \psi(\mathbf{x}, t) d\mathbf{x}, \quad t \geq 0. \quad (2.4)$$

Other very useful quantities characterizing the dynamics of rotating BEC in 2D are the condensate width defined as

$$\delta_x(t) = \int_{\mathbb{R}^d} x^2 |\psi(\mathbf{x}, t)|^2 d\mathbf{x}, \quad \delta_y(t) = \int_{\mathbb{R}^d} y^2 |\psi(\mathbf{x}, t)|^2 d\mathbf{x}, \quad \delta_r(t) = \delta_x(t) + \delta_y(t). \quad (2.5)$$

As proven in [5], in 2D with a radial symmetric trap, i.e.,  $d = 2$  and  $\gamma_x = \gamma_y := \gamma_r$  in (1.3), for any initial data  $\psi_0(x, y)$  in (1.4), we have for any  $t \geq 0$

$$\delta_r(t) = \frac{E_{\beta,\Omega}(\psi_0) + \Omega \langle L_z \rangle(0)}{\gamma_r^2} [1 - \cos(2\gamma_r t)] + \delta_r^{(0)} \cos(2\gamma_r t) + \frac{\delta_r^{(1)}}{2\gamma_r} \sin(2\gamma_r t), \quad (2.6)$$

where  $\delta_r^{(0)} := \delta_r(0) = \delta_x(0) + \delta_y(0)$  and  $\delta_r^{(1)} := \dot{\delta}_r(0) = \dot{\delta}_x(0) + \dot{\delta}_y(0)$  with for  $\alpha = x$  or  $y$

$$\begin{aligned} \delta_\alpha(0) &= \delta_\alpha^{(0)} = \int_{\mathbb{R}^2} \alpha^2 |\psi_0(\mathbf{x})|^2 d\mathbf{x}, \\ \dot{\delta}_\alpha(0) &= \delta_\alpha^{(1)} = 2 \int_{\mathbb{R}^2} \alpha \left[ -\Omega |\psi_0|^2 (x \partial_y - y \partial_x) \alpha + \text{Im}(\psi_0^* \partial_\alpha \psi_0) \right] d\mathbf{x}. \end{aligned}$$

Furthermore, when the initial data  $\psi_0(x, y)$  in (1.4) satisfies

$$\psi_0(x, y) = f(r) e^{im\theta} \quad \text{with} \quad m \in \mathbb{Z} \quad \text{and} \quad f(0) = 0 \quad \text{when} \quad m \neq 0, \quad (2.7)$$

with  $(r, \theta)$  being the polar coordinates in 2D, we have, for any  $t \geq 0$ ,

$$\begin{aligned} \delta_x(t) &= \delta_y(t) = \frac{1}{2} \delta_r(t) \\ &= \frac{E_{\beta,\Omega}(\psi_0) + m\Omega}{2\gamma_x^2} [1 - \cos(2\gamma_x t)] + \delta_x^{(0)} \cos(2\gamma_x t) + \frac{\delta_x^{(1)}}{2\gamma_x} \sin(2\gamma_x t). \end{aligned} \quad (2.8)$$

These immediately imply that  $\delta_r(t)$  is a periodic function with angular frequency doubling the trapping frequency in 2D with a radial symmetric trap, and also  $\delta_x(t)$  and  $\delta_y(t)$  are periodic functions with frequency doubling the trapping frequency provided that the initial data satisfies (2.7).

### 3 A time-splitting pseudospectral method for rotating BEC

In this section, we will present an explicit, unconditionally stable and spectrally accurate numerical method to solve the GPE (1.3) for dynamics of rotating BEC.

Due to the external trapping potential  $V_d(\mathbf{x})$ , the solution  $\psi(\mathbf{x}, t)$  of (1.3)-(1.4) decays to zero exponentially fast when  $|\mathbf{x}| \rightarrow \infty$ . Thus in practical computation, we always truncate the problem (1.3)-(1.4) into a bounded computational domain with homogeneous Dirichlet boundary condition:

$$i \partial_t \psi(\mathbf{x}, t) = -\frac{1}{2} \nabla^2 \psi + [V(\mathbf{x}) - iW(\mathbf{x})] \psi + \beta_d |\psi|^2 \psi - \Omega L_z \psi, \quad \mathbf{x} \in \Omega_{\mathbf{x}}, \quad t > 0, \quad (3.1)$$

$$\psi(\mathbf{x}, t) = 0, \quad \mathbf{x} \in \Gamma = \partial\Omega_{\mathbf{x}}, \quad t \geq 0, \quad (3.2)$$

$$\psi(\mathbf{x}, 0) = \psi_0(\mathbf{x}), \quad \mathbf{x} \in \bar{\Omega}_{\mathbf{x}}; \quad (3.3)$$

where  $W(\mathbf{x}) \geq 0$  corresponds to a localized loss term [36] and  $V(\mathbf{x}) = V_d(\mathbf{x}) + V_p(\mathbf{x})$  with  $V_p(\mathbf{x}) \geq 0$  a conservative repulsive pinning potential [36]. Here we choose  $\Omega_{\mathbf{x}} = [a, b] \times [c, d]$  in 2D, and resp.,  $\Omega_{\mathbf{x}} = [a, b] \times [c, d] \times [e, f]$  in 3D, with  $|a|$ ,  $b$ ,  $|c|$ ,  $d$ ,  $|e|$  and  $f$  sufficiently large. The use of more sophisticated radiation boundary conditions is an interesting topic that remains to be examined in the future.

#### 3.1 Time-splitting

We choose a time step size  $\Delta t > 0$ . For  $n = 0, 1, 2, \dots$ , from time  $t = t_n = n\Delta t$  to  $t = t_{n+1} = t_n + \Delta t$ , the GPE (3.1) is first solved in two splitting steps. One solves first

$$i \partial_t \psi(\mathbf{x}, t) = -\frac{1}{2} \nabla^2 \psi(\mathbf{x}, t) - \Omega L_z \psi(\mathbf{x}, t) \quad (3.4)$$

for the time step of length  $\Delta t$ , followed by solving

$$i \partial_t \psi(\mathbf{x}, t) = [V(\mathbf{x}) - iW(\mathbf{x})] \psi(\mathbf{x}, t) + \beta_d |\psi(\mathbf{x}, t)|^2 \psi(\mathbf{x}, t), \quad (3.5)$$

for the same time step. For  $t \in [t_n, t_{n+1}]$ , multiplying (3.5) by  $\psi^*$ , the conjugate of  $\psi$ , we get

$$i \psi^*(\mathbf{x}, t) \partial_t \psi(\mathbf{x}, t) = [V(\mathbf{x}) - iW(\mathbf{x})] |\psi(\mathbf{x}, t)|^2 + \beta_d |\psi(\mathbf{x}, t)|^4. \quad (3.6)$$

Subtracting the conjugate of (3.6) from (3.6) and multiplying by  $-i$ , one obtains

$$\frac{d}{dt} |\psi(\mathbf{x}, t)|^2 = \psi^* \partial_t \psi + \psi \partial_t \psi^* = -2W(\mathbf{x}) |\psi(\mathbf{x}, t)|^2. \quad (3.7)$$

Solving (3.7), we get

$$|\psi(\mathbf{x}, t)|^2 = e^{-2W(\mathbf{x})(t-t_n)} |\psi(\mathbf{x}, t_n)|^2, \quad t_n \leq t \leq t_{n+1}. \quad (3.8)$$

Substituting (3.8) into (3.5), we obtain

$$i \partial_t \psi(\mathbf{x}, t) = [V(\mathbf{x}) - iW(\mathbf{x})] \psi(\mathbf{x}, t) + \beta_d e^{-2W(\mathbf{x})(t-t_n)} |\psi(\mathbf{x}, t_n)|^2 \psi(\mathbf{x}, t). \quad (3.9)$$

Integrate (3.9) from  $t_n$  to  $t$ , we find for  $\mathbf{x} \in \Omega_{\mathbf{x}}$  and  $t_n \leq t \leq t_{n+1}$ :

$$\psi(\mathbf{x}, t) = \begin{cases} e^{-i[V(\mathbf{x}) + \beta_d |\psi(\mathbf{x}, t_n)|^2](t-t_n)} \psi(\mathbf{x}, t_n), & W(\mathbf{x}) = 0, \\ \frac{\psi(\mathbf{x}, t_n)}{e^{W(\mathbf{x})(t-t_n)}} e^{-i[V(\mathbf{x})(t-t_n) + \beta_d |\psi(\mathbf{x}, t_n)|^2(1-e^{-2W(\mathbf{x})(t-t_n)})/2W(\mathbf{x})]}, & W(\mathbf{x}) > 0. \end{cases} \quad (3.10)$$

To discretize (3.4) in space, compared with non-rotating BEC [7, 11, 6], i.e.  $\Omega = 0$  in (1.3), the main difficulty is that the coefficients in  $L_z$  are *not* constants which cause big trouble in applying sine or Fourier pseudospectral discretization. Due to the special structure in the angular momentum rotation term  $L_z$  (1.5), we will apply the alternating direction implicit (ADI) technique and decouple the operator  $-\frac{1}{2}\nabla^2 - \Omega L_z$  into two one dimensional operators such that each operator becomes a summation of terms with constant coefficients in that dimension. Therefore, they can be discretized in space by Fourier pseudospectral method and the ODEs in phase space can be integrated analytically. The details for discretizing (3.4) in 2D & 3D will be presented in the next two subsections respectively.

## 3.2 Discretization in 2D

When  $d = 2$  in (3.4), we choose mesh sizes  $\Delta x > 0$  and  $\Delta y > 0$  with  $\Delta x = (b-a)/M$  and  $\Delta y = (d-c)/N$  for  $M$  and  $N$  even positive integers, and let the grid points be

$$x_j = a + j\Delta x, \quad j = 0, 1, 2, \dots, M; \quad y_k = c + k\Delta y, \quad k = 0, 1, 2, \dots, N.$$

Let  $\psi_{jk}^n$  be the approximation of  $\psi(x_j, y_k, t_n)$  and  $\psi^n$  be the solution vector with component  $\psi_{jk}^n$ .

From time  $t = t_n$  to  $t = t_{n+1}$ , we solve (3.4) first

$$i \partial_t \psi(\mathbf{x}, t) = -\frac{1}{2} \partial_{xx} \psi(\mathbf{x}, t) - i\Omega y \partial_x \psi(\mathbf{x}, t), \quad (3.11)$$

for the time step of length  $\Delta t$ , followed by solving

$$i \partial_t \psi(\mathbf{x}, t) = -\frac{1}{2} \partial_{yy} \psi(\mathbf{x}, t) + i\Omega x \partial_y \psi(\mathbf{x}, t), \quad (3.12)$$

for the same time step. The detailed discretizations of (3.11) and (3.12) are shown in Appendix A.

## 3.3 Discretization in 3D

When  $d = 3$  in (3.4), we choose mesh sizes  $\Delta x > 0$ ,  $\Delta y > 0$  and  $\Delta z > 0$  with  $\Delta x = (b-a)/M$ ,  $\Delta y = (d-c)/N$  and  $\Delta z = (f-e)/L$  for  $M$ ,  $N$  and  $L$  even positive integers, and let the grid points be

$$x_j = a + j\Delta x, \quad 0 \leq j \leq M; \quad y_k = c + k\Delta y, \quad 0 \leq k \leq N; \quad z_l = e + l\Delta z, \quad 0 \leq l \leq L.$$

Let  $\psi_{jkl}^n$  be the approximation of  $\psi(x_j, y_k, z_l, t_n)$  and  $\psi^n$  be the solution vector with component  $\psi_{jkl}^n$ .

Similar as those for 2D case, from time  $t = t_n$  to  $t = t_{n+1}$ , we solve (3.4) first

$$i \partial_t \psi(\mathbf{x}, t) = \left( -\frac{1}{2} \partial_{xx} - \frac{1}{4} \partial_{zz} - i \Omega y \partial_x \right) \psi(\mathbf{x}, t), \quad (3.13)$$

for the time step of length  $\Delta t$ , followed by solving

$$i \partial_t \psi(\mathbf{x}, t) = \left( -\frac{1}{2} \partial_{yy} - \frac{1}{4} \partial_{zz} + i \Omega x \partial_y \right) \psi(\mathbf{x}, t), \quad (3.14)$$

for the same time step. The detailed discretizations of (3.13) and (3.14) are shown in Appendix B.

### 3.4 Stability

We define the usual discrete  $l^2$ -norm for the solution  $\psi^n$  as

$$\|\psi^n\|_{l^2} = \sqrt{\frac{b-a}{M} \frac{d-c}{N} \sum_{j=0}^{M-1} \sum_{k=0}^{N-1} |\psi_{jk}^n|^2}, \quad (3.15)$$

for  $d = 2$ , and for  $d = 3$

$$\|\psi^n\|_{l^2} = \sqrt{\frac{b-a}{M} \frac{d-c}{N} \frac{f-e}{L} \sum_{j=0}^{M-1} \sum_{k=0}^{N-1} \sum_{l=0}^{L-1} |\psi_{jkl}^n|^2}. \quad (3.16)$$

For the *stability* of the time-splitting spectral approximations (A.7) for 2D and (B.1) for 3D, we have the following lemma, which shows that the total density is conserved when  $W(\mathbf{x}) \equiv 0$ , and resp. decreased when  $W(\mathbf{x}) > 0$ , in the discretized level.

**Lemma 3.1** *The time-splitting spectral schemes (A.7) for 2D and (B.1) for 3D GPE with an angular momentum rotation term are unconditionally stable. In fact, for any mesh sizes  $\Delta x > 0$ ,  $\Delta y > 0$  and  $\Delta z > 0$ , and time step size  $\Delta t > 0$ ,*

$$\|\psi^n\|_{l^2} \leq \|\psi^{n-1}\|_{l^2} \leq \|\psi^0\|_{l^2} = \|\psi_0\|_{l^2}, \quad n = 1, 2, \dots. \quad (3.17)$$

*In addition, if  $W(\mathbf{x}) \equiv 0$  in (3.1), then we have*

$$\|\psi^n\|_{l^2} \equiv \|\psi^0\|_{l^2} = \|\psi_0\|_{l^2}, \quad n = 1, 2, \dots. \quad (3.18)$$

**Proof:** Follows the line of the analogous results for the linear and nonlinear Schrödinger equations in [6, 9, 10, 7, 12].

## 4 Numerical results

In this section, we first test the accuracy of our new numerical method (A.7) for 2D and (B.1) for 3D and compare our numerical results with the analytical results reviewed in section 2. Then we apply our new numerical method to study vortex lattice dynamics in rotating BEC by changing the trapping frequencies and to generate a giant vortex by introducing a localized loss term. Our aim is not to find new physics phenomena but to demonstrate the efficiency and high resolution of our new numerical method

### 4.1 Accuracy test

To test the accuracy of our numerical method in 2D, we take  $\beta_2 = 100$  and  $\Omega = 0.7$  in (1.3). The initial condition in (1.4) is taken as

$$\psi_0(x, y) = \frac{(\gamma_x \gamma_y)^{\frac{1}{4}}}{\pi^{\frac{1}{2}}} e^{-(\gamma_x x^2 + \gamma_y y^2)/2}, \quad \mathbf{x} = (x, y)^T \in \mathbb{R}^2. \quad (4.1)$$

We take  $\gamma_x = 1.0$  and  $\gamma_y = 2.0$  in (1.3) and (4.1). Similar example was used in [5, 41] for testing numerical accuracy of different numerical methods for rotating BEC. The GPE (1.3) is solved on  $[-8, 8] \times [-8, 8]$ , i.e. we take  $a = -8$ ,  $b = 8$ ,  $c = -8$  and  $d = 8$ . Let  $\psi$  be the *exact* solution which is obtained numerically by using our method with a very fine mesh and small time step, e.g.  $\Delta x = \Delta y = \frac{1}{64}$  and  $\Delta t = 0.0001$ , and  $\psi^{(\Delta x, \Delta y, \Delta t)}$  be the numerical solution obtained with the mesh size  $(\Delta x, \Delta y)$  and time step  $\Delta t$ .

First we test the spectral accuracy in space by choosing a very small time step  $\Delta t = 0.0001$ , and solving the problem for each fixed  $\beta_2$  with different mesh size  $\Delta x = \Delta y$  so that the discretization errors in time can be neglected comparing to those in space. The errors  $\|\psi(t) - \psi^{(\Delta x, \Delta y, \Delta t)}(t)\|_{l^2}$  at  $t = 0.5$  are shown in Table 1 for different values  $\beta_2$  and  $h = \Delta x = \Delta y$ .

h	1/2	1/4	1/8
$\beta_2 = 20$	2.68E-2	6.5E-5	3.41E-10
$\beta_2 = 50$	0.1315	2.01E-3	5.91E-8
$\beta_2 = 100$	0.4287	1.94E-2	8.89E-6

Table 1: Spatial discretization errors  $\|\psi(t) - \psi^{(\Delta x, \Delta y, \Delta t)}(t)\|_{l^2}$  at  $t = 0.5$  in 2D.

Next we test the second-order accuracy in time. Table 2 lists the errors at  $t = 0.5$  for different values of  $\beta_2$  and time steps  $\Delta t$  with a fine mesh in space, i.e.  $\Delta x = \Delta y = 1/16$ .

Similarly, to test the accuracy of our numerical method in 3D, we take  $\beta_3 = 100$



$\Delta t$	1/40	1/80	1/160	1/320	1/640
$\beta_2 = 20$	7.86E-4	1.95E-4	4.86E-5	1.21E-5	3.02E-6
$\beta_2 = 50$	2.95E-3	7.24E-4	1.80E-4	4.49E-5	1.12E-5
$\beta_2 = 100$	7.98E-3	1.98E-3	4.76E-4	1.19E-4	2.96E-5

Table 2: Temporal discretization errors  $\|\psi(t) - \psi^{(\Delta x, \Delta y, \Delta t)}(t)\|_{l^2}$  at  $t = 0.5$  in 2D.

and  $\Omega = 0.7$  in (1.3). The initial condition in (1.4) is taken as

$$\psi_0(x, y, z) = \frac{(\gamma_x \gamma_y \gamma_z)^{\frac{1}{4}}}{\pi^{\frac{3}{4}}} e^{-(\gamma_x x^2 + \gamma_y y^2 + \gamma_z z^2)/2}, \quad \mathbf{x} = (x, y, z)^T \in \mathbb{R}^3. \quad (4.2)$$

We take  $\gamma_x = \gamma_y = \gamma_z = 1.0$  in (1.3) and (4.2), and solve the GPE (1.3) in 3D on  $[-8, 8] \times [-8, 8] \times [-8, 8]$ . Again let  $\psi$  be the *exact* solution which is obtained numerically by using our method with a fine mesh and small time step, e.g.  $\Delta x = \Delta y = \Delta z = \frac{1}{8}$  and  $\Delta t = 0.0001$ , and  $\psi^{(\Delta x, \Delta y, \Delta z, \Delta t)}$  be the numerical solution obtained with mesh size  $(\Delta x, \Delta y, \Delta z)$  and time step  $\Delta t$ . Table 3 shows the spatial discretization errors  $\|\psi(t) - \psi^{(\Delta x, \Delta y, \Delta z, \Delta t)}(t)\|_{l^2}$  at  $t = 0.5$  with  $\Delta t = 0.0001$  for different values  $\beta_3$  and mesh sizes  $h = \Delta x = \Delta y = \Delta z$ . Table 4 lists the errors at  $t = 0.5$  for different values of  $\beta_3$  and time steps  $\Delta t$  with a fine mesh in space, i.e.  $\Delta x = \Delta y = \Delta z = 1/8$ .

h	1	1/2	1/4
$\beta_3 = 20$	5.78E-2	1.27E-3	2.38E-8
$\beta_3 = 50$	0.1515	9.50E-3	2.45E-6
$\beta_3 = 100$	0.3075	3.88E-2	7.08E-5

Table 3: Spatial discretization errors  $\|\psi(t) - \psi^{(\Delta x, \Delta y, \Delta, \Delta z, t)}(t)\|_{l^2}$  at  $t = 0.5$  in 3D.

$\Delta t$	1/40	1/80	1/160	1/320	1/640
$\beta_3 = 20$	1.778E-4	4.435E-5	1.108E-5	2.767E-6	6.897E-7
$\beta_3 = 50$	6.266E-4	1.559E-4	3.892E-5	9.718E-6	2.422E-6
$\beta_3 = 100$	1.63E-3	4.0379E-4	1.0069E-4	2.5141E-5	6.265E-6

Table 4: Temporal discretization errors  $\|\psi(t) - \psi^{(\Delta x, \Delta y, \Delta z, \Delta t)}(t)\|_{l^2}$  at  $t = 0.5$  in 3D.

From Tables 1-4, we can draw the following conclusions: (i) The method (A.7) or (B.1) is of spectral accuracy in space and second order accuracy in time. (ii) For a given fixed mesh size and time step, when  $\beta_d$  is increasing, the errors are increasing too. This implies that when the number of atoms in the condensate is increasing,

i.e.  $\beta_d$  is increasing, more grid points and smaller time step are needed in practical computation in order to achieve a given accuracy.

Furthermore, Fig. 1 shows time evolution of the normalization  $N(\psi)(t)$ , energy  $E_{\beta,\Omega}(\psi)(t)$ , angular momentum expectation  $\langle L_z \rangle(t)$  and condensate widths for the above parameters setup in 2D and 3D.

From Fig. 1, we can see that: (i). The normalization is conserved in both cases which confirms (2.1). (ii). The energy is not conserved in the discretized level, but the perturbation is very small, e.g. less than 5% (c.f. (a), (c), (e) and (g) of Fig. 1). (iii). The angular momentum expectation is conserved when  $\gamma_x = \gamma_y = 1$  (c.f. (a) and (e) of Fig. 1) which confirms the analytical result (2.3), and oscillates when  $1 = \gamma_x \neq \gamma_y = 2$  (c.f. (c) and (g) of Fig. 1). (iv) The condensate widths  $\delta_x(t)$ ,  $\delta_y(t)$  and  $\delta_r(t)$  are periodic functions when  $\gamma_x = \gamma_y = 1$  which confirm the analytical result (2.6) (c.f. (b) of Fig. 1), and are not periodic functions when  $1 = \gamma_x \neq \gamma_y = 2$  (c.f. (d) and (h) of Fig. 1).

## 4.2 Dynamics of a vortex lattice in rotating BEC

In this subsection we numerically study the dynamics of a vortex lattice in rotating BEC by changing trap frequencies. This study was motivated by the recent experiment [20] in which the frequencies of trapping potential of a stable BEC were changed [20]. One of the most striking observation in the experiment is that the condensate contains sheet-like structures rather than individual vortex cores in the dynamics by deforming the static trap [20]. By using the hydrodynamic forms of the GPE (1.3) in Thomas-Fermi regime, Cozzin et al. [18] tried to study this phenomena theoretically. But they did not find the sheet-like structures by changing the trap frequencies in their theoretical study. Here we study this phenomena by directly simulating the GPE (1.3) using our new numerical method.

We take  $d = 2$ ,  $\beta_2 = 100$  and  $\Omega = 0.99$  in (1.3). The initial data  $\psi_0(\mathbf{x})$  in (1.4) is chosen as the ground state of (1.3) with  $d = 2$ ,  $\Omega = 0.99$ ,  $\beta_2 = 100$  and  $\gamma_x = \gamma_y = 1$ , which is computed numerically by the normalized gradient flow proposed in [13]. In the ground state, there are about 61 vortices in the vortex lattice (c.f. Fig. 2). We solve the problem on  $\Omega_{\mathbf{x}} = [-24, 24] \times [-24, 24]$  with mesh size  $\Delta x = \Delta y = 3/16$  and time step  $\Delta t = 0.001$ .

First we study free expansion of the quantized vortex lattice. In general, the size of a stable vortex lattice in a BEC is too small to visualize it. In experiments, by removing the trap, i. e., letting the vortex lattice freely expands, one can obtain an enlarged vortex lattice so as to take a photo for it [20]. Of course, they hope the vortex structure doesn't change during the free expansion. Thus theoretical study of free expansion is very helpful for experiments. We start with the stable BEC and remove the trapping at time  $t = 0$ , i.e. choosing  $\gamma_x = \gamma_y = 0$  in (1.3). Similar numerical study was also carried out in [2] by a different numerical method with much less number of vortices in the lattice. Fig. 2 shows image plots of the density  $|\psi(\mathbf{x}, t)|^2$  at different times for the free expansion of the vortex lattice. From

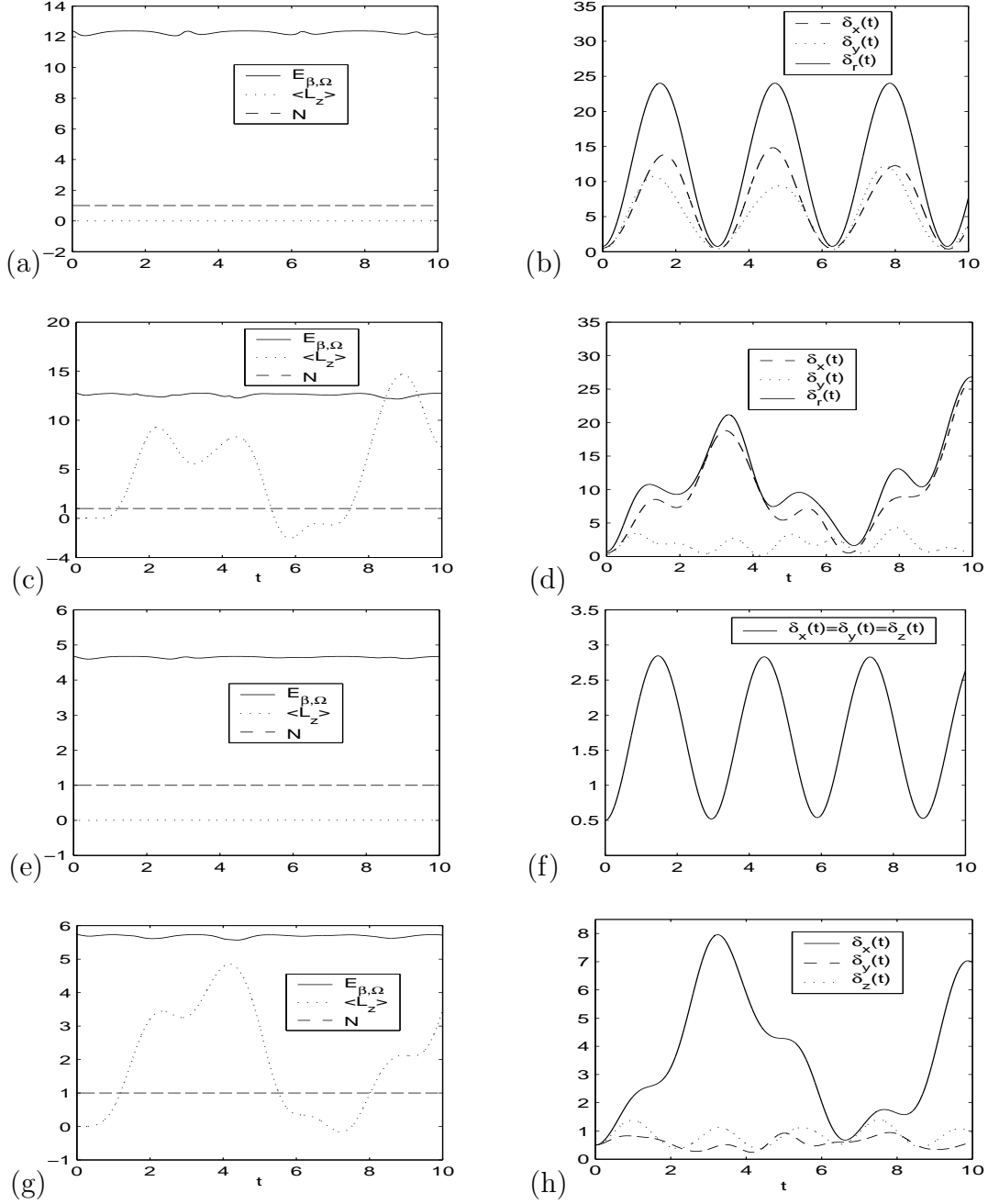


Fig. 1: Time evolution of the normalization  $N(t) := N(\psi)(t)$ , energy  $E_{\beta,\Omega}(\psi)(t)$ , angular momentum expectation  $\langle L_z \rangle(t)$  (left column), and condensate widths  $\delta_x(t)$ ,  $\delta_y(t)$  and  $\delta_r(t)$  (right column). Results in 2D for  $\gamma_x = \gamma_y = 1$  (a&b), and  $\gamma_x = 1$  and  $\gamma_y = 2$  (c&d); and results in 3D for  $\gamma_x = \gamma_y = \gamma_z = 1$  (e&f), and  $\gamma_x = 1$ ,  $\gamma_y = 2$  and  $\gamma_z = 1.5$  (g&h).

the figure, we can see that when the trap is removed at  $t = 0$ , the vortex lattice will expand with time and the vortex structure as well as the rotational symmetry is kept during the expansion. This gives a numerical justification for the free expansion

used in BEC experiments.

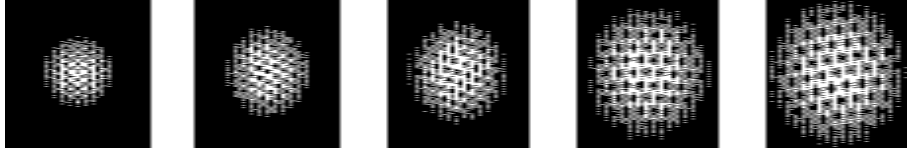


Fig. 2: Image plots of the density  $|\psi(\mathbf{x}, t)|^2$  on  $[-18, 18] \times [-18, 18]$  at different times  $t = 0, 0.75, 1.5, 2.0$  and  $2.75$  (from left to right) for the free expansion of a quantized vortex lattice.

Next, we study dynamics of the quantized vortex lattice by changing the trap frequencies. We study six different cases: I.  $\gamma_x = 1, \gamma_y = 1.5$ ; II.  $\gamma_x = 1, \gamma_y = 0.75$ ; III.  $\gamma_x = 1.5, \gamma_y = 1$ ; IV.  $\gamma_x = 0.75, \gamma_y = 1$ ; V.  $\gamma_x = \sqrt{1.2}, \gamma_y = \sqrt{0.8}$ ; VI.  $\gamma_x = \sqrt{1.4}, \gamma_y = \sqrt{0.6}$ . Similar numerical study was also carried out in [2] by a different method with much less number of vortices in the lattice. Fig. 3 shows image plots of the density  $|\psi(\mathbf{x}, t)|^2$  at different times for cases I-II for changing frequencies in  $y$ -direction only. Fig. 4 shows similar results for cases III-IV for changing frequencies in  $x$ -direction only, and Fig. 5 for cases V-VI for changing frequencies in both  $x$ - and  $y$ -directions.

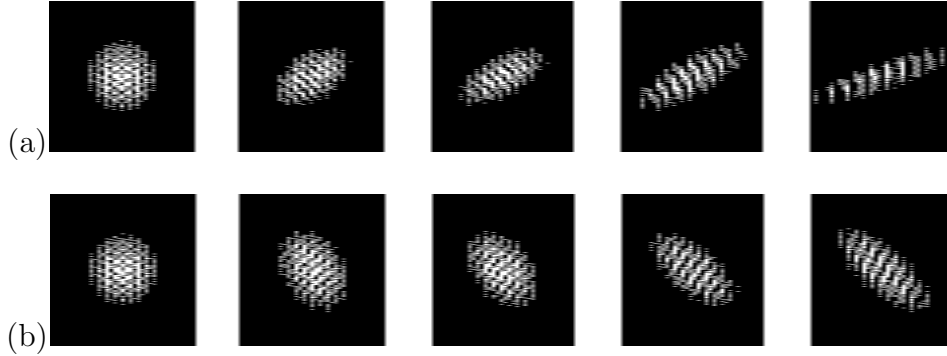


Fig. 3: Image plots of the density  $|\psi(\mathbf{x}, t)|^2$  on  $[-18, 18] \times [-18, 18]$  at different times  $t=0, 2.0, 4.0, 6.0$ , and  $8.0$  (from left to right) by changing frequency in  $y$ -direction only from  $\gamma_y = 1$  to: (a)  $\gamma_y = 1.5$ ; (b)  $\gamma_y = 0.75$ .

In Figs. 3-5, initially the condensate is assumed to be in its ground state which is a vortex lattice with about 61 vortices. From the numerical results presented here, when the trap frequencies are changed at  $t = 0$ , we find that: (i) Cases I & II correspond to changing trap frequency in  $y$ -direction only. The condensate initially starts to contract (c.f. (a) of Fig. 3) or expand (c.f. (b) of Fig. 3) in  $y$ -direction since the trap frequency in  $y$ -direction is increasing or decreasing at  $t = 0$ . (ii) Cases III & IV correspond to changing trap frequency in  $x$ -direction only. Similar results are observed (c.f. (a) and (b) of Fig. 4). (iii) Cases V & VI correspond to increasing and decreasing the trapping frequencies in  $x$  and  $y$ -directions by the same value, i.e.,  $\epsilon$ , respectively [18]. The condensate initially starts to contract and expand in  $x$  and  $y$ -directions respectively (c.f. Fig. 5). (iv) We numerically observed

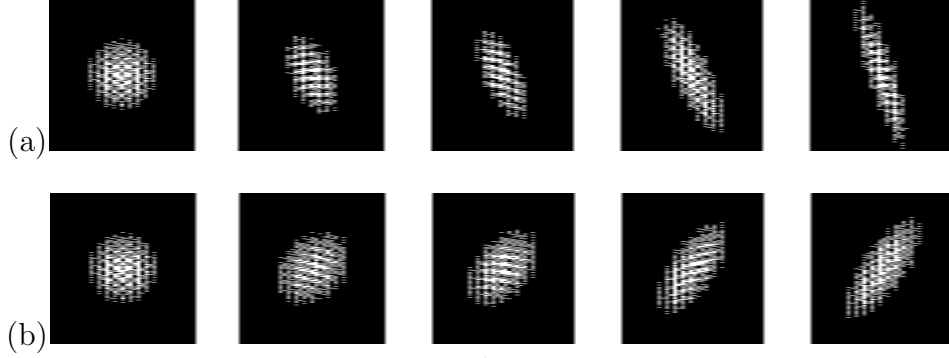


Fig. 4: Image plots of the density  $|\psi(\mathbf{x}, t)|^2$  on  $[-18, 18] \times [-18, 18]$  at different times  $t=0, 2.0, 4.0, 6.0$ , and  $8.0$  (from left to right) by changing frequency in  $x$ -direction only from  $\gamma_x = 1$  to: (a)  $\gamma_x = 1.5$ ; (b)  $\gamma_x = 0.75$ .

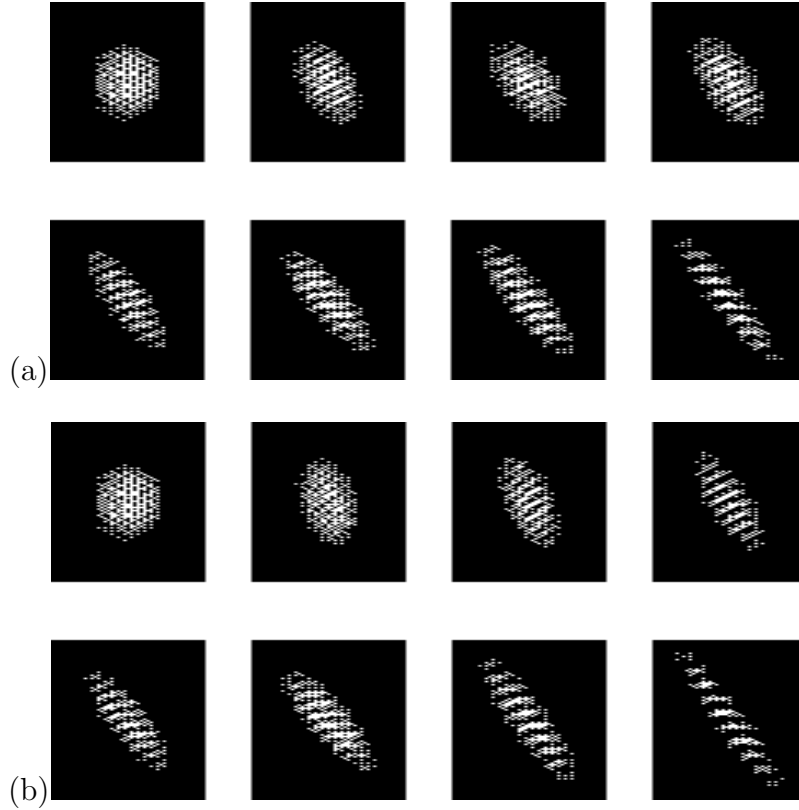


Fig. 5: Image plots of the density  $|\psi(\mathbf{x}, t)|^2$  on  $[-18, 18] \times [-18, 18]$  at different times by changing trapping frequencies from  $\gamma_x = 1, \gamma_y = 1$  to  $\gamma_x = \sqrt{1 + \epsilon}, \gamma_y = \sqrt{1.0 - \epsilon}$ . (a) With  $\epsilon = 0.2$  for times  $t = 0, 2, 3, 4, 5, 6, 7$  and  $8$  (from left to right); (b) with  $\epsilon = 0.4$  for times  $t = 0, 1.0, 1.5, 2.0, 2.5, 3.0, 4.0$  and  $4.5$ .

the remarkable sheet-like vortices in our numerical results (c.f. Figs. 3&5). One can compare our numerical results in Figs. 3&5 with the experimental results, e.g. Fig. 4 in [20], and find very good qualitative agreement in sheet-like vortex lattice formation. Furthermore, we also found that when  $\epsilon$  is bigger in Cases V&VI, the

sheet-like vortices appear earlier.

### 4.3 Generation of giant vortex in rotating BEC

In this subsection we numerically generate a giant vortex in rotating BEC from its ground state by introducing a localized loss term [21]. This study was motivated by the recent experiment [21] and theoretical study [36] in which the giant vortex formation arises as a dynamics effect. We take  $d = 2$ ,  $\beta_2 = 100$ ,  $V_p(\mathbf{x}) \equiv 0$  and  $\Omega = 0.99$  in (3.1). The initial data  $\psi_0(\mathbf{x})$  in (3.3) is chosen as the ground state of (1.3) with  $d = 2$ ,  $\Omega = 0.99$ ,  $\beta_2 = 100$  and  $\gamma_x = \gamma_y = 1$ , which is computed numerically by the normalized gradient flow proposed in [13]. The localized loss term in (3.1) is chosen as a Gaussian function of the form [36]

$$W(x, y) = w_0 \exp \left[ -\frac{(x - x_0)^2 + (y - y_0)^2}{r_0^2} \right], \quad (x, y) \in \mathbb{R}^2, \quad (4.3)$$

where  $w_0$ ,  $x_0$ ,  $y_0$  and  $r_0$  are constants. We take  $w_0 = 1$  and  $r_0 = \sqrt{7}/2$  in (4.3) and solve the problem (3.1)-(3.3) on  $\Omega_{\mathbf{x}} = [-24, 24] \times [-24, 24]$  with mesh size  $\Delta x = \Delta y = 3/16$  and time step  $\Delta t = 0.001$ . Fig. 6 shows image plots of the density  $|\psi(\mathbf{x}, t)|^2$  at different times for different values of  $(x_0, y_0)$ .

From Fig. 6, we can see that the giant vortex lattice is generated due to the dynamic effect in a rotating BEC. The center of the giant vortex is the same as the center of the localized loss term and the size of the giant vortex depends on the values of  $r_0$  and  $w_0$ . One can compare our numerical results in Fig. 6 with the experimental results, e.g. Fig. 1 in [21], and the theoretical study, e.g. Fig. 1 in [36], and find very good qualitative agreement in giant-vortex formation.

## 5 Conclusions

We have proposed a new time-splitting Fourier pseudospectral method for computing dynamics of rotating BEC based on an efficient approximation of GPE with an angular momentum rotation term. The new method is explicit, unconditionally stable, and of spectral accuracy in space and second order accuracy in time. It is time reversible and time transverse invariant in the discretized level, just as the original GPE does, and conserves the total density in the discretized level. The efficient and accurate numerical method was applied to study dynamics of a quantized vortex lattice in rotating BEC. In the future, we plan to extend the idea for constructing the new numerical method for one component rotating BEC to multi-component rotating BEC and spinor dynamics in a rotational frame, and apply the method to study vortex line dynamics in rotating BEC in 3D.

## Appendix A. Discretization in 2D

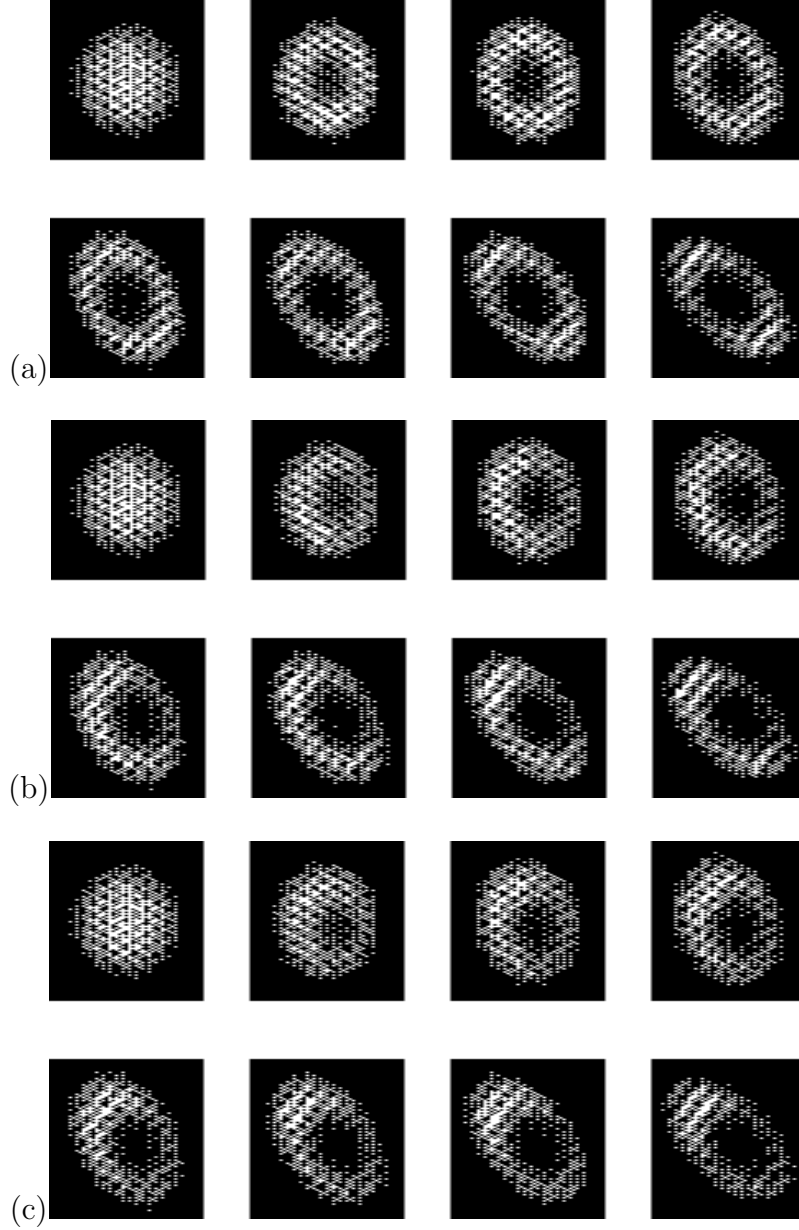


Fig. 6: Image plots of the density  $|\psi(\mathbf{x}, t)|^2$  on  $[-12, 12] \times [-12, 12]$  at different times  $t=0, 0.5, 0.75, 1.0, 1.25, 1.5, 1.75$ , and  $2.0$  (from left to right) with  $w_0 = 1$  and  $r_0 = \frac{\sqrt{7}}{2}$  in (4.3) for generating giant vortices. (a)  $x_0 = 0, y_0 = 0$ ; (b)  $x_0 = 1.5, y_0 = 0$ ; (c)  $x_0 = 1.5, y_0 = 1$ .

For each fixed  $y$ , the operator in the equation (3.11) is in  $x$ -direction with constant coefficients and thus we can discretize it in  $x$ -direction by a Fourier pseudospectral method. Assume

$$\psi(x, y, t) = \sum_{p=-M/2}^{M/2-1} \hat{\psi}_p(y, t) \exp[i\mu_p(x - a)], \quad (\text{A.1})$$

where  $\mu_p = \frac{2p\pi}{b-a}$  and  $\widehat{\psi}_p(y, t)$  is the Fourier coefficient for the  $p$ -th mode in  $x$ -direction. Plugging (A.1) into (3.11), noticing the orthogonality of the Fourier functions, we obtain for  $-\frac{M}{2} \leq p \leq \frac{M}{2} - 1$  and  $c \leq y \leq d$ :

$$i \partial_t \widehat{\psi}_p(y, t) = \left( \frac{1}{2} \mu_p^2 + \Omega y \mu_p \right) \widehat{\psi}_p(y, t), \quad t_n \leq t \leq t_{n+1}. \quad (\text{A.2})$$

The above linear ODE can be integrated in time *exactly* and we obtain

$$\widehat{\psi}_p(y, t) = \exp \left[ -i \left( \frac{1}{2} \mu_p^2 + \Omega y \mu_p \right) (t - t_n) \right] \widehat{\psi}_p(y, t_n), \quad t_n \leq t \leq t_{n+1}. \quad (\text{A.3})$$

Similarly, for each fixed  $x$ , the operator in the equation (3.12) is in  $y$ -direction with constant coefficients and thus we can discretize it in  $y$ -direction by a Fourier pseudospectral method. Assume

$$\psi(x, y, t) = \sum_{q=-N/2}^{N/2-1} \widehat{\psi}_q(x, t) \exp[i\lambda_q(y - c)], \quad (\text{A.4})$$

where  $\lambda_q = \frac{2q\pi}{d-c}$  and  $\widehat{\psi}_q(x, t)$  is the Fourier coefficient for the  $q$ -th mode in  $y$ -direction. Plugging (A.4) into (3.12), noticing the orthogonality of the Fourier functions, we obtain for  $-\frac{N}{2} \leq q \leq \frac{N}{2} - 1$  and  $a \leq x \leq b$ :

$$i \partial_t \widehat{\psi}_q(x, t) = \left( \frac{1}{2} \lambda_q^2 - \Omega x \lambda_q \right) \widehat{\psi}_q(x, t), \quad t_n \leq t \leq t_{n+1}. \quad (\text{A.5})$$

Again the above linear ODE can be integrated in time *exactly* and we obtain

$$\widehat{\psi}_q(x, t) = \exp \left[ -i \left( \frac{1}{2} \lambda_q^2 - \Omega x \lambda_q \right) (t - t_n) \right] \widehat{\psi}_q(x, t_n), \quad t_n \leq t \leq t_{n+1}. \quad (\text{A.6})$$

From time  $t = t_n$  to  $t = t_{n+1}$ , we combine the splitting steps via the standard second order Strang splitting [37, 4, 12]:

$$\begin{aligned} \psi_{jk}^{(1)} &= \sum_{p=-M/2}^{M/2-1} e^{-i\Delta t(\mu_p^2 + 2\Omega y_k \mu_p)/4} \widehat{(\psi_k^n)}_p e^{i\mu_p(x_j - a)}, \quad 0 \leq j \leq M, \quad 0 \leq k \leq N, \\ \psi_{jk}^{(2)} &= \sum_{q=-N/2}^{N/2-1} e^{-i\Delta t(\lambda_q^2 - 2\Omega x_j \lambda_q)/4} \widehat{(\psi_j^{(1)})}_q e^{i\lambda_q(y_k - c)}, \quad 0 \leq k \leq N, \quad 0 \leq j \leq M, \\ \psi_{jk}^{(3)} &= \begin{cases} e^{-i\Delta t[V(x_j, y_k) + \beta_2 |\psi_{jk}^{(2)}|^2]} \psi_{jk}^{(2)}, & W(x_j, y_k) = 0, \\ \frac{\psi_{jk}^{(2)}}{e^{\Delta t W(x_j, y_k)}} e^{-i[\Delta t V(x_j, y_k) + \beta_2 |\psi_{jk}^{(2)}|^2 (1 - e^{-2\Delta t W(x_j, y_k)})/2W(x_j, y_k)]}, & W(x_j, y_k) > 0, \end{cases} \\ \psi_{jk}^{(4)} &= \sum_{q=-N/2}^{N/2-1} e^{-i\Delta t(\lambda_q^2 - 2\Omega x_j \lambda_q)/4} \widehat{(\psi_j^{(3)})}_q e^{i\lambda_q(y_k - c)}, \quad 0 \leq k \leq N, \quad 0 \leq j \leq M, \\ \psi_{jk}^{n+1} &= \sum_{p=-M/2}^{M/2-1} e^{-i\Delta t(\mu_p^2 + 2\Omega y_k \mu_p)/4} \widehat{(\psi_k^{(4)})}_p e^{i\mu_p(x_j - a)}, \quad 0 \leq j \leq M, \quad 0 \leq k \leq N; \quad (\text{A.7}) \end{aligned}$$



where for each fixed  $k$ ,  $(\widehat{\psi_k^\alpha})_p$  ( $p = -M/2, \dots, M/2 - 1$ ) with  $\alpha$  an index, the Fourier coefficients of the vector  $\psi_k^\alpha = (\psi_{0k}^\alpha, \psi_{1k}^\alpha, \dots, \psi_{(M-1)k}^\alpha)^T$ , are defined as

$$(\widehat{\psi_k^\alpha})_p = \frac{1}{M} \sum_{j=0}^{M-1} \psi_{jk}^\alpha e^{-i\mu_p(x_j-a)}, \quad p = -\frac{M}{2}, \dots, \frac{M}{2} - 1; \quad (\text{A.8})$$

similarly, for each fixed  $j$ ,  $(\widehat{\psi_j^\alpha})_q$  ( $q = -N/2, \dots, N/2 - 1$ ), the Fourier coefficients of the vector  $\psi_j^\alpha = (\psi_{j0}^\alpha, \psi_{j1}^\alpha, \dots, \psi_{j(N-1)}^\alpha)^T$ , are defined as

$$(\widehat{\psi_j^\alpha})_q = \frac{1}{N} \sum_{k=0}^{N-1} \psi_{jk}^\alpha e^{-i\lambda_q(y_k-c)}, \quad q = -\frac{N}{2}, \dots, \frac{N}{2} - 1. \quad (\text{A.9})$$

For the algorithm (A.7) presented in Appendix A, the total memory requirement is  $O(MN)$  and the total computational cost per time step is  $O(MN \ln(MN))$ . The scheme is time reversible when  $W(\mathbf{x}) \equiv 0$ , just as it holds for the GPE (1.3), i.e. the scheme is unchanged if we interchange  $n \leftrightarrow n+1$  and  $\Delta t \leftrightarrow -\Delta t$  in (A.7). Also, a main advantage of the numerical method is its time-transverse invariance when  $W(\mathbf{x}) \equiv 0$ , just as it holds for the GPE (1.3) itself. If a constant  $\alpha$  is added to the external potential  $V$ , then the discrete wave functions  $\psi_{jk}^{n+1}$  obtained from (A.7) get multiplied by the phase factor  $e^{-i\alpha(n+1)\Delta t}$ , which leaves the discrete quadratic observable  $|\psi_{jk}^{n+1}|^2$  unchanged. This property does not hold for the finite difference scheme [25, 39], the leap-frog spectral method [41] and the efficient discretization proposed in [5] for GPE with an angular momentum term.

## Appendix B. Discretization in 3D

For each fixed  $y$ , the operator in the equation (3.13) is in  $x$  and  $z$ -directions with constant coefficients and thus we can discretize it in  $x$  and  $z$ -directions by a Fourier pseudospectral method. Similarly, for each fixed  $x$ , the operator in the equation (3.14) is in  $y$  and  $z$ -directions with constant coefficients and thus we can discretize it in  $y$  and  $z$ -directions by a Fourier pseudospectral method. The discretizations of (3.13) and (3.14) are similar as those for (3.11) and (3.12) respectively and they are omitted here. For simplicity and convenience of the reader, here we only present the algorithm for 3D GPE with an angular momentum rotation term (1.3) with  $0 \leq j \leq M$ ,  $0 \leq k \leq N$  and  $0 \leq l \leq L$ :

$$\begin{aligned} \psi_{jkl}^{(1)} &= \sum_{p=-M/2}^{M/2-1} \sum_{s=-L/2}^{L/2-1} e^{-i\Delta t(2\mu_p^2 + \gamma_s^2 + 4\Omega y_k \mu_p)/8} (\widehat{\psi_k^n})_{ps} e^{i\mu_p(x_j-a)} e^{i\gamma_s(z_l-e)}, \\ \psi_{jkl}^{(2)} &= \sum_{q=-N/2}^{N/2-1} \sum_{s=-L/2}^{L/2-1} e^{-i\Delta t(2\lambda_q^2 + \gamma_s^2 - 4\Omega x_j \lambda_q)/8} (\widehat{\psi_j^{(1)}})_{qs} e^{i\lambda_q(y_k-c)} e^{i\gamma_s(z_l-e)}, \end{aligned}$$

$$\begin{aligned}
\psi_{jkl}^{(3)} &= \begin{cases} e^{-i\Delta t[V(x_j, y_k, z_l) + \beta_3 |\psi_{jkl}^{(2)}|^2]} \psi_{jkl}^{(2)}, & W(x_j, y_k, z_l) = 0, \\ \frac{\psi_{jkl}^{(2)}}{e^{\Delta t W(x_j, y_k, z_l)}} \exp[-i(\Delta t V(x_j, y_k, z_l) + \beta_3 |\psi_{jkl}^{(2)}|^2 (1 - e^{-2\Delta t W(x_j, y_k, z_l)})/2W(x_j, y_k, z_l))], & W(x_j, y_k, z_l) > 0, \end{cases} \\
\psi_{jkl}^{(4)} &= \sum_{q=-N/2}^{N/2-1} \sum_{s=-L/2}^{L/2-1} e^{-i\Delta t(2\lambda_q^2 + \gamma_s^2 - 4\Omega x_j \lambda_q)/8} \widehat{(\psi_j^{(3)})}_{qs} e^{i\lambda_q(y_k - c)} e^{i\gamma_s(z_l - e)}, \\
\psi_{jkl}^{n+1} &= \sum_{p=-M/2}^{M/2-1} \sum_{s=-L/2}^{L/2-1} e^{-i\Delta t(2\mu_p^2 + \gamma_s^2 + 4\Omega y_k \mu_p)/8} \widehat{(\psi_k^{(4)})}_{ps} e^{i\mu_p(x_j - a)} e^{i\gamma_s(z_l - e)}, \quad (\text{B.1})
\end{aligned}$$

where for each fixed  $k$ ,  $\widehat{(\psi_k^\alpha)}_{ps}$  ( $-M/2 \leq p \leq M/2 - 1$ ,  $-L/2 \leq s \leq L/2 - 1$ ) with  $\alpha$  an index, the Fourier coefficients of the vector  $\psi_{jkl}^\alpha$  ( $0 \leq j < M$ ,  $0 \leq l < L$ ), are defined as

$$\widehat{(\psi_k^\alpha)}_{ps} = \frac{1}{ML} \sum_{j=0}^{M-1} \sum_{l=0}^{L-1} \psi_{jkl}^\alpha e^{-i\mu_p(x_j - a)} e^{-i\gamma_s(z_l - e)}, \quad -\frac{M}{2} \leq p < \frac{M}{2}, \quad -\frac{L}{2} \leq s < \frac{L}{2};$$

similarly, for each fixed  $j$ ,  $\widehat{(\psi_j^\alpha)}_{qs}$  ( $-N/2 \leq q \leq N/2 - 1$ ,  $-L/2 \leq s \leq L/2 - 1$ ) with  $\alpha$  an index, the Fourier coefficients of the vector  $\psi_{jkl}^\alpha$  ( $k = 0, \dots, N$ ,  $l = 0, \dots, L$ ), are defined as

$$\widehat{(\psi_j^\alpha)}_{qs} = \frac{1}{NL} \sum_{m=0}^{N-1} \sum_{l=0}^{L-1} \psi_{jlm}^\alpha e^{-i\lambda_q(y_k - c)} e^{-i\gamma_s(z_l - e)}, \quad -\frac{N}{2} \leq q < \frac{N}{2}, \quad -\frac{L}{2} \leq s < \frac{L}{2};$$

with  $\gamma_s = \frac{2\pi s}{f - e}$  for  $s = -L/2, \dots, L/2 - 1$ .

For the discretization in 3D, the total memory requirement is  $O(MNL)$  and the total computational cost per time step is  $O(MNL \ln(MNL))$ . Furthermore, the discretization is time reversible and time transverse invariant in the discretized level when  $W(\mathbf{x}) \equiv 0$ .

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$\beta_3=100.0$

